

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF051917\
 Data File : BF095272.D
 Acq On : 19 May 2017 20:02
 Operator : SJ/MA
 Sample : I3203-06MSD
 Misc :
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 E2-T8-BASEMSD

Manual Integrations
 APPROVED

mohammad
 5/22/2017 1:29:49 PM

Quant Time: May 22 07:45:19 2017
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF050217.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu May 18 17:07:53 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.77	152	80965	20.00	ng	0.00
21) Naphthalene-d8	9.78	136	363678	20.00	ng	0.00
38) Acenaphthene-d10	12.61	164	178574	20.00	ng	0.00
63) Phenanthrene-d10	15.02	188	322541	20.00	ng	0.00
75) Chrysene-d12	18.69	240	246051	20.00	ng	0.00
86) Perylene-d12	20.37	264	166338	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.74	112	699921	144.13	ng	0.00
7) Phenol-d6	7.31	99	859805	147.56	ng	0.00
23) Nitrobenzene-d5	8.67	82	548849	95.17	ng	0.00
41) 2,4,6-Tribromophenol	13.92	330	220358	150.68	ng	0.00
44) 2-Fluorobiphenyl	11.55	172	1006009	81.09	ng	0.00
78) Terphenyl-d14	17.34	244	858211	76.82	ng	-0.01

Target Compounds						Qvalue
2) 1,4-Dioxane	2.27	88	90188	37.87	ng	92
3) Pyridine	2.98	79	182442	33.05	ng	95
4) n-Nitrosodimethylamine	2.90	42	81611	35.47	ng	91
6) Aniline	7.28	93	285870	38.86	ng	95
8) 2-Chlorophenol	7.46	128	273907	49.55	ng	94
9) Benzaldehyde	7.10	77	85334	23.98	ng	97
10) Phenol	7.33	94	286750	46.70	ng	91
11) bis(2-Chloroethyl)ether	7.41	93	252851	49.88	ng	94
12) 1,3-Dichlorobenzene	7.67	146	282798	45.10	ng	99
13) 1,4-Dichlorobenzene	7.80	146	290337	45.66	ng	97
14) 1,2-Dichlorobenzene	8.02	146	285478	48.10	ng	98
15) Benzyl Alcohol	8.05	79	224193	59.82	ng	96
16) 2,2'-oxybis(1-Chloropropan	8.25	45	339424	47.31	ng	57
17) 2-Methylphenol	8.26	107	231429	51.16	ng	# 91
18) Hexachloroethane	8.54	117	100059	46.45	ng	# 85
19) n-Nitroso-di-n-propylamine	8.47	70	205742	52.21	ng	92
20) 3+4-Methylphenols	8.52	107	313935	51.07	ng	# 60
22) Acetophenone	8.44	105	406143	46.55	ng	# 84
24) Nitrobenzene	8.70	77	283297	46.86	ng	90
25) Isophorone	9.10	82	525701	51.05	ng	95
26) 2-Nitrophenol	9.21	139	167101	51.23	ng	97
27) 2,4-Dimethylphenol	9.34	122	264075	50.07	ng	98
28) bis(2-Chloroethoxy)methane	9.46	93	351940	49.40	ng	100
29) 2,4-Dichlorophenol	9.62	162	246543	49.51	ng	98
30) 1,2,4-Trichlorobenzene	9.71	180	242448	45.45	ng	98
31) Naphthalene	9.81	128	840177	45.97	ng	100
32) Benzoic acid	9.62	122	55357	19.38	ng	98
33) 4-Chloroaniline	9.99	127	77318	11.35	ng	94
34) Hexachlorobutadiene	10.03	225	117691	41.81	ng	97
35) Caprolactam	10.60	113	71426m	47.77	ng	
36) 4-Chloro-3-methylphenol	10.81	107	285220	53.57	ng	88
37) 2-Methylnaphthalene	10.94	142	600385	50.05	ng	98
39) 1,2,4,5-Tetrachlorobenzene	11.22	216	235824	42.94	ng	99
40) Hexachlorocyclopentadiene	11.19	237	116053	52.52	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	11.44	196	172921	47.16	ng	98
43) 2,4,5-Trichlorophenol	11.54	196	162407	41.21	ng	92
45) 1,1'-Biphenyl	11.71	154	665746	44.28	ng	98
46) 2-Chloronaphthalene	11.72	162	532156	43.83	ng	95
47) 2-Nitroaniline	11.95	65	155286	46.73	ng	# 78
48) Acenaphthylene	12.38	152	842934	44.33	ng	98
49) Dimethylphthalate	12.26	163	852436	58.15	ng	99
50) 2,6-Dinitrotoluene	12.35	165	141443	47.29	ng	93
51) Acenaphthene	12.67	154	458472	40.37	ng	98
52) 3-Nitroaniline	12.64	138	61384	19.12	ng	89
53) 2,4-Dinitrophenol	12.84	184	83048	52.73	ng	# 66
54) Dibenzofuran	12.95	168	731249	46.53	ng	98
55) 4-Nitrophenol	13.02	139	156989	81.43	ng	# 72
56) 2,4-Dinitrotoluene	13.02	165	191439	49.81	ng	# 92
57) Fluorene	13.51	166	621259	45.46	ng	99
58) 2,3,4,6-Tetrachlorophenol	13.20	232	129452	52.19	ng	96
59) Diethylphthalate	13.41	149	638937	45.47	ng	100
60) 4-Chlorophenyl-phenylether	13.53	204	276543	46.04	ng	90
61) 4-Nitroaniline	13.63	138	123644	41.96	ng	97
62) Azobenzene	13.79	77	585903	51.89	ng	94
64) 4,6-Dinitro-2-methylphenol	13.69	198	95183	44.02	ng	# 70
65) n-Nitrosodiphenylamine	13.74	169	537989	45.96	ng	98
66) 4-Bromophenyl-phenylether	14.32	248	155945	45.76	ng	98
67) Hexachlorobenzene	14.41	284	157457	45.09	ng	# 88
68) Atrazine	14.67	200	162161	49.06	ng	95
69) Pentachlorophenol	14.77	266	129065	81.61	ng	99
70) Phenanthrene	15.06	178	839808	45.32	ng	98
71) Anthracene	15.14	178	883549	46.67	ng	99
72) Carbazole	15.42	167	835935	48.53	ng	97
73) Di-n-butylphthalate	15.98	149	1050320	48.90	ng	99
74) Fluoranthene	16.80	202	853122	44.88	ng	99
76) Benzidine	17.02	184	265918	41.06	ng	# 93
77) Pyrene	17.10	202	861902	46.84	ng	99
79) Butylbenzylphthalate	17.99	149	438446	51.22	ng	# 85
80) Benzo(a)anthracene	18.68	228	665027	46.44	ng	99
81) 3,3'-Dichlorobenzidine	18.66	252	147264	29.78	ng	# 97
82) Chrysene	18.72	228	625181	45.15	ng	98
83) Bis(2-ethylhexyl)phthalate	18.74	149	617537	50.64	ng	# 100
84) Di-n-octyl phthalate	19.51	149	994547	49.74	ng	96
85) Indeno(1,2,3-cd)pyrene	21.69	276	463851	41.25	ng	99
87) Benzo(b)fluoranthene	19.95	252	466675	45.40	ng	98
88) Benzo(k)fluoranthene	19.98	252	551853m	55.66	ng	
89) Benzo(a)pyrene	20.31	252	455716	48.05	ng	99
90) Dibenzo(a,h)anthracene	21.69	278	378834	48.36	ng	98
91) Benzo(g,h,i)perylene	22.08	276	410165	53.36	ng	98

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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